

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_Y\Data\VY071624\  
 Data File : VY018842.D  
 Acq On : 16 Jul 2024 20:01  
 Operator : SY/MD  
 Sample : VSTDCCC050  
 Misc : 5.01g/5.0mL/MSVOA\_Y/SOIL/A  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 MSVOA\_Y  
 LabSampleId :  
 VSTDCCC050

Quant Time: Jul 17 02:01:07 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_Y\methods\82Y070924S.M  
 Quant Title : SW846 8260  
 QLast Update : Tue Jul 09 14:28:51 2024  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|       | Compound                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I   | Pentafluorobenzene          | 1.000 | 1.000 | 0.0    | 83    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.419 | 0.409 | 2.4    | 75    | 0.00     |
| 3 P   | Chloromethane               | 0.637 | 0.648 | -1.7   | 82    | 0.00     |
| 4 C   | Vinyl Chloride              | 0.745 | 0.764 | -2.6#  | 81    | 0.00     |
| 5 T   | Bromomethane                | 0.512 | 0.557 | -8.8   | 95    | 0.00     |
| 6 T   | Chloroethane                | 0.484 | 0.504 | -4.1   | 87    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.973 | 1.004 | -3.2   | 80    | 0.00     |
| 8 T   | Diethyl Ether               | 0.237 | 0.277 | -16.9  | 95    | 0.00     |
| 9 T   | 1,1,2-Trichlorotrifluoroeth | 0.499 | 0.499 | 0.0    | 77    | 0.00     |
| 10 T  | Methyl Iodide               | 0.538 | 0.529 | 1.7    | 73    | 0.00     |
| 11 T  | Tert butyl alcohol          | 0.050 | 0.040 | 20.0   | 89    | 0.00     |
| 12 CM | 1,1-Dichloroethene          | 0.483 | 0.481 | 0.4#   | 80    | 0.00     |
| 13 T  | Acrolein                    | 0.040 | 0.024 | 40.0#  | 55    | 0.00     |
| 14 T  | Allyl chloride              | 0.598 | 0.625 | -4.5   | 85    | 0.00     |
| 15 T  | Acrylonitrile               | 0.094 | 0.115 | -22.3  | 100   | 0.00     |
| 16 T  | Acetone                     | 0.091 | 0.088 | 3.3    | 82    | 0.00     |
| 17 T  | Carbon Disulfide            | 1.353 | 1.344 | 0.7    | 78    | 0.00     |
| 18 T  | Methyl Acetate              | 0.264 | 0.423 | -60.2# | 143   | 0.00     |
| 19 T  | Methyl tert-butyl Ether     | 1.121 | 1.321 | -17.8  | 95    | 0.00     |
| 20 T  | Methylene Chloride          | 0.610 | 0.628 | -3.0   | 97    | -0.01    |
| 21 T  | trans-1,2-Dichloroethene    | 0.510 | 0.538 | -5.5   | 84    | 0.00     |
| 22 T  | Diisopropyl ether           | 1.302 | 1.419 | -9.0   | 91    | 0.00     |
| 23 T  | Vinyl Acetate               | 1.026 | 1.073 | -4.6   | 83    | 0.00     |
| 24 P  | 1,1-Dichloroethane          | 0.831 | 0.887 | -6.7   | 88    | 0.00     |
| 25 T  | 2-Butanone                  | 0.127 | 0.145 | -14.2  | 94    | 0.00     |
| 26 T  | 2,2-Dichloropropane         | 0.771 | 0.676 | 12.3   | 72    | 0.00     |
| 27 T  | cis-1,2-Dichloroethene      | 0.594 | 0.642 | -8.1   | 90    | 0.00     |
| 28 T  | Bromochloromethane          | 0.334 | 0.383 | -14.7  | 95    | 0.00     |
| 29 T  | Tetrahydrofuran             | 0.078 | 0.095 | -21.8  | 97    | 0.00     |
| 30 C  | Chloroform                  | 0.926 | 1.051 | -13.5# | 94    | 0.00     |
| 31 T  | Cyclohexane                 | 0.756 | 0.700 | 7.4    | 77    | 0.00     |
| 32 T  | 1,1,1-Trichloroethane       | 0.861 | 0.913 | -6.0   | 84    | 0.00     |
| 33 S  | 1,2-Dichloroethane-d4       | 0.492 | 0.559 | -13.6  | 89    | 0.00     |
| 34 I  | 1,4-Difluorobenzene         | 1.000 | 1.000 | 0.0    | 89    | 0.00     |
| 35 S  | Dibromofluoromethane        | 0.313 | 0.329 | -5.1   | 85    | 0.00     |
| 36 T  | 1,1-Dichloropropene         | 0.432 | 0.417 | 3.5    | 81    | 0.00     |
| 37 T  | Ethyl Acetate               | 0.180 | 0.200 | -11.1  | 91    | 0.00     |
| 38 T  | Carbon Tetrachloride        | 0.521 | 0.516 | 1.0    | 81    | 0.00     |
| 39 T  | Methylcyclohexane           | 0.554 | 0.541 | 2.3    | 80    | 0.00     |
| 40 TM | Benzene                     | 1.254 | 1.296 | -3.3   | 89    | 0.00     |
| 41 T  | Methacrylonitrile           | 0.102 | 0.106 | -3.9   | 95    | -0.03    |
| 42 TM | 1,2-Dichloroethane          | 0.351 | 0.380 | -8.3   | 93    | 0.00     |
| 43 T  | Isopropyl Acetate           | 0.357 | 0.404 | -13.2  | 96    | 0.00     |
| 44 TM | Trichloroethene             | 0.358 | 0.376 | -5.0   | 91    | 0.00     |
| 45 C  | 1,2-Dichloropropane         | 0.273 | 0.286 | -4.8#  | 90    | 0.00     |
| 46 T  | Dibromomethane              | 0.177 | 0.203 | -14.7  | 96    | 0.00     |
| 47 T  | Bromodichloromethane        | 0.442 | 0.487 | -10.2  | 94    | 0.00     |
| 48 T  | Methyl methacrylate         | 0.162 | 0.184 | -13.6  | 95    | 0.00     |

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 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 MSVOA\_Y  
 LabSampleId :  
 VSTDCCC050

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 Quant Title : SW846 8260  
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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|       | Compound                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|--------|-------|----------|
| 49 T  | 1,4-Dioxane                 | 0.002 | 0.002 | 0.0    | 99    | 0.00     |
| 50 S  | Toluene-d8                  | 1.186 | 1.209 | -1.9   | 82    | 0.00     |
| 51 T  | 4-Methyl-2-Pentanone        | 0.182 | 0.219 | -20.3  | 99    | 0.00     |
| 52 CM | Toluene                     | 0.833 | 0.861 | -3.4#  | 89    | 0.00     |
| 53 T  | t-1,3-Dichloropropene       | 0.382 | 0.415 | -8.6   | 93    | 0.00     |
| 54 T  | cis-1,3-Dichloropropene     | 0.448 | 0.478 | -6.7   | 90    | 0.00     |
| 55 T  | 1,1,2-Trichloroethane       | 0.224 | 0.252 | -12.5  | 97    | 0.00     |
| 56 T  | Ethyl methacrylate          | 0.295 | 0.342 | -15.9  | 97    | 0.00     |
| 57 T  | 1,3-Dichloropropane         | 0.371 | 0.418 | -12.7  | 96    | 0.00     |
| 58 T  | 2-Chloroethyl Vinyl ether   | 0.141 | 0.179 | -27.0# | 100   | 0.00     |
| 59 T  | 2-Hexanone                  | 0.128 | 0.149 | -16.4  | 96    | 0.00     |
| 60 T  | Dibromochloromethane        | 0.308 | 0.356 | -15.6  | 98    | 0.00     |
| 61 T  | 1,2-Dibromoethane           | 0.218 | 0.255 | -17.0  | 97    | 0.00     |
| 62 S  | 4-Bromofluorobenzene        | 0.405 | 0.437 | -7.9   | 89    | 0.00     |
| 63 I  | Chlorobenzene-d5            | 1.000 | 1.000 | 0.0    | 94    | 0.00     |
| 64 T  | Tetrachloroethene           | 0.405 | 0.458 | -13.1  | 99    | 0.00     |
| 65 PM | Chlorobenzene               | 1.055 | 1.068 | -1.2   | 91    | 0.00     |
| 66 T  | 1,1,1,2-Tetrachloroethane   | 0.363 | 0.375 | -3.3   | 93    | 0.00     |
| 67 C  | Ethyl Benzene               | 1.848 | 1.791 | 3.1#   | 87    | 0.00     |
| 68 T  | m/p-Xylenes                 | 0.725 | 0.712 | 1.8    | 88    | 0.00     |
| 69 T  | o-Xylene                    | 0.682 | 0.675 | 1.0    | 88    | 0.00     |
| 70 T  | Styrene                     | 1.137 | 1.178 | -3.6   | 92    | 0.00     |
| 71 P  | Bromoform                   | 0.198 | 0.233 | -17.7  | 103   | 0.00     |
| 72 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 98    | 0.00     |
| 73 T  | Isopropylbenzene            | 3.659 | 3.407 | 6.9    | 86    | 0.00     |
| 74 T  | N-amyl acetate              | 0.748 | 0.779 | -4.1   | 91    | 0.00     |
| 75 P  | 1,1,2,2-Tetrachloroethane   | 0.577 | 0.578 | -0.2   | 93    | 0.00     |
| 76 T  | 1,2,3-Trichloropropane      | 0.392 | 0.422 | -7.7   | 108   | 0.00     |
| 77 T  | Bromobenzene                | 0.874 | 0.874 | 0.0    | 94    | 0.00     |
| 78 T  | n-propylbenzene             | 4.346 | 3.952 | 9.1    | 84    | 0.00     |
| 79 T  | 2-Chlorotoluene             | 2.443 | 2.253 | 7.8    | 87    | 0.00     |
| 80 T  | 1,3,5-Trimethylbenzene      | 2.994 | 2.843 | 5.0    | 88    | 0.00     |
| 81 T  | trans-1,4-Dichloro-2-butene | 0.175 | 0.177 | -1.1   | 89    | 0.00     |
| 82 T  | 4-Chlorotoluene             | 2.516 | 2.348 | 6.7    | 88    | 0.00     |
| 83 T  | tert-Butylbenzene           | 2.764 | 2.571 | 7.0    | 87    | 0.00     |
| 84 T  | 1,2,4-Trimethylbenzene      | 2.975 | 2.825 | 5.0    | 89    | 0.00     |
| 85 T  | sec-Butylbenzene            | 3.924 | 3.634 | 7.4    | 84    | 0.00     |
| 86 T  | p-Isopropyltoluene          | 3.387 | 3.083 | 9.0    | 84    | 0.00     |
| 87 T  | 1,3-Dichlorobenzene         | 1.716 | 1.681 | 2.0    | 92    | 0.00     |
| 88 T  | 1,4-Dichlorobenzene         | 1.701 | 1.620 | 4.8    | 90    | 0.00     |
| 89 T  | n-Butylbenzene              | 3.071 | 2.796 | 9.0    | 82    | 0.00     |
| 90 T  | Hexachloroethane            | 0.630 | 0.592 | 6.0    | 86    | 0.00     |
| 91 T  | 1,2-Dichlorobenzene         | 1.482 | 1.501 | -1.3   | 96    | 0.00     |
| 92 T  | 1,2-Dibromo-3-Chloropropane | 0.093 | 0.108 | -16.1  | 99    | 0.00     |
| 93 T  | 1,2,4-Trichlorobenzene      | 0.905 | 0.944 | -4.3   | 93    | 0.00     |
| 94 T  | Hexachlorobutadiene         | 0.521 | 0.484 | 7.1    | 81    | 0.00     |
| 95 T  | Naphthalene                 | 1.552 | 1.824 | -17.5  | 102   | 0.00     |

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 96 T | 1,2,3-Trichlorobenzene | 0.774 | 0.825 | -6.6 | 95    | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 6